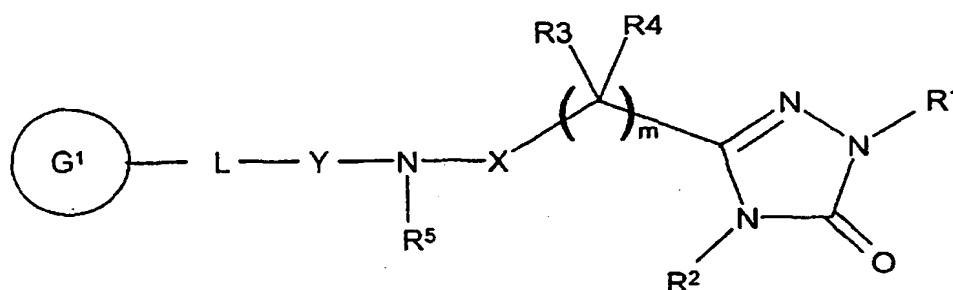


CLAIMS

1. A compound of formula (I) or a pharmaceutically acceptable salt or solvate thereof



(I)

wherein

R^1 and R^2 independently represent H or C1 to 6 alkyl; said alkyl being optionally further substituted by an aryl ring or an aromatic heterocyclic ring containing 1 to 3 heteroatoms independently selected from O, S and N; said aromatic ring being optionally further substituted by halogen, CF_3 , C1 to 4 alkyl or C1 to 4 alkoxy;

Each R^3 and each R^4 independently represents H or C1 to 6 alkyl; said alkyl being optionally further substituted by OH, C1 to 4 alkoxy, C1 to 4 alkylthio, amino, N-alkylamino or N,N-dialkylamino;

or R^3 and R^4 are bonded together so as to form a 3 to 7 membered ring; said ring optionally incorporating one heteroatom selected from O, $S(O)_q$ and N;

m represents an integer 1, 2 or 3;

X represents a group S(O), S(O)₂ or C(=O);

R⁵ represents H or C1 to 6 alkyl; said alkyl being optionally further substituted by halogen,
5 OH or C1 to 6 alkoxy;

Y represents a direct bond;

or Y and R⁵ are bonded together such that the group -NR⁵Y- together represents a 4 to 7
10 membered saturated or partially unsaturated azacyclic ring; said azacyclic ring optionally
incorporating one further heteroatom selected from O, S(O)_n and N; said azacyclic ring
being optionally benzo fused; said azacyclic ring being optionally substituted by C1 to 6
alkyl, C1 to 6 alkoxy or OH;

15 L represents a direct bond;

or L represents O, S(O)_p, C(O), NR⁶, C(O)NR⁶, NR⁶C(O), C2 to 6 alkynyl, C2 to 6
alkenyl, C1 to 6 alkyl, C1 to 6 heteroalkyl or C3 to 6 heteroalkynyl; said alkyl, alkenyl or
alkynyl group being optionally further substituted by halogen, OH or C1 to 6 alkoxy;

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n, p and q independently represent an integer 0, 1 or 2;

G¹ represents a monocyclic, bicyclic, tricyclic or tetracyclic group comprising one, two,
three or four ring structures each of up to 7 ring atoms; each ring structure being
25 independently selected from cycloalkyl; cycloalkenyl; heterocycloalkyl; unsaturated
heterocycloalkyl; aryl; or an aromatic heterocyclic ring containing 1 to 3 heteroatoms
independently selected from O, S and N; with each ring structure being independently
optionally substituted by one or more substituents independently selected from halogen,
hydroxy, CHO, C1 to 6 alkyl, C1 to 6 alkoxy, halo-C1 to 6 alkoxy, amino, N-alkylamino,

N,N-dialkylamino, alkylsulfonamino, C2 to 6 alkanoylamino, cyano, nitro, thiol, alkylthio, alkylsulfonyl, alkylaminosulfonyl, C2 to 6 alkanoyl, aminocarbonyl, N-alkylamino-carbonyl, N,N-amino-carbonyl;

- 5 wherein any alkyl radical within any substituent may itself be optionally substituted with one or more groups selected from halogen, hydroxy, C1 to 6 alkoxy, halo-C1 to 6 alkoxy, amino, N-alkylamino, N,N-dialkylamino, N-alkylsulfonamino, N-C2 to 6 alkanoylamino, cyano, nitro, thiol, alkylthio, alkylsulfonyl, N-alkylaminosulfonyl, CHO, C2 to 6 alkanoyl, aminocarbonyl, N-alkylaminocarbonyl, N,N-dialkylaminocarbonyl and carbamate;

10 and wherein any alkyl radical is a C1 to 6 alkyl radical;

- and when G^1 is a bicyclic, tricyclic or tetracyclic group, each ring structure is independently joined to the next ring structure by a direct bond, by -O-, by C1-6 alkyl, by C1-6 haloalkyl, by C1-6 heteroalkyl, by C2-6 alkenyl, by C2-6 alkynyl, by sulfone, by CO, by NR^7CO , by $CONR^7$, by NR^7 , by S, or by C(OH), or each ring structure is fused to the next ring structure;

- 20 R^6 and R^7 independently represent H or C1 to 6 alkyl;

and when the group $-NR^5Y-$ represents an azacyclic ring and L represents a direct bond, the group G^1 may also be spiro fused to the azacyclic ring;

- 25 2. A compound according to claim 1, wherein X represents $S(O)_2$.

3. A compound according to claim 1 or 2, wherein R^1 and R^2 each represent hydrogen.

4. A compound according to any one of claims 1 to 3, wherein R^3 and R^4 each represent hydrogen.

5. A compound according to any one of claims 1 to 4, wherein R^5 represents hydrogen or C1 to 6 alkyl and Y represents a direct bond.

6. A compound according to any one of claims 1 to 4, wherein the group $-NR^5Y-$ together represents a five or six membered saturated or partially unsaturated azacyclic ring, said azacyclic ring optionally incorporating one further heteroatom selected from O, $S(O)_n$ and N.

7. A compound according to any one of claims 1 to 6 wherein L represents a direct bond, O, C2 to 6 alkynyl, C1 to 6 alkyl, C1 to 6 heteroalkyl or C3 to 6 heteroalkynyl.

8. A compound according to any one of claims 1 to 7, wherein G^1 represents an optionally substituted monocyclic or bicyclic ring structure.

9. A compound according to claim 1 which is selected from the group consisting of:

5-[(4-[(5-chloropyridin-2-yl)oxy]piperidin-1-yl)sulfonyl)methyl]-2,4-dihydro-3H-1,2,4-triazol-3-one;

5-[2-[(4-[(5-chloropyridin-2-yl)oxy]piperidin-1-yl)sulfonyl]ethyl]-2,4-dihydro-3H-1,2,4-triazol-3-one;

5-[3-[(4-[(5-chloropyridin-2-yl)oxy]piperidin-1-yl)sulfonyl]propyl]-2,4-dihydro-3H-1,2,4-triazol-3-one;

5-[(4-(4-chlorophenyl)piperazin-1-yl)sulfonyl)methyl]-2,4-dihydro-3H-1,2,4-triazol-3-one;

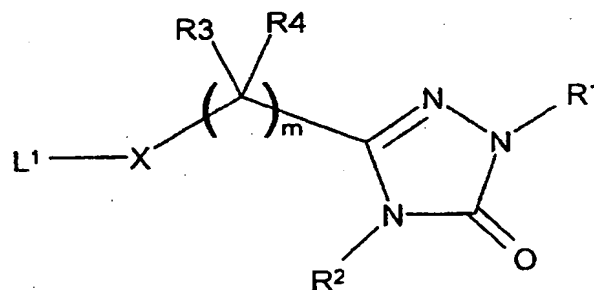
5-[(4-[(2-methoxypyrimidin-5-yl)ethynyl]-3,6-dihydropyridin-1(2H)-yl)sulfonyl)methyl]-2,4-dihydro-3H-1,2,4-triazol-3-one;

5-[(4-[(2-(trifluoromethyl)pyrimidin-5-yl)ethynyl]-3,6-dihydropyridin-1(2H)-yl)sulfonyl)methyl]-2,4-dihydro-3H-1,2,4-triazol-3-one;

- 5-([4-[(2-cyclopropyl)pyrimidin-5-yl]ethynyl]-3,6-dihydropyridin-1(2H)-yl)sulfonyl)methyl)-2,4-dihydro-3H-1,2,4-triazol-3-one;
 5-([4-(4-chlorophenyl)piperidin-1-yl]sulfonyl)methyl)-2,4-dihydro-3H-1,2,4-triazol-3-one;
 5 N-benzyl-1-(5-oxo-4,5-dihydro-1H-1,2,4-triazol-3-yl)methanesulfonamide;
 1-(5-oxo-4,5-dihydro-1H-1,2,4-triazol-3-yl)-N-(2-phenylethyl)methanesulfonamide;
 5-(2-{[4-(4-chlorophenyl)piperidin-1-yl]sulfonyl}ethyl)-2,4-dihydro-3H-1,2,4-triazol-3-one;
 5-(2-{[4-(4-chlorophenyl)piperazin-1-yl]sulfonyl}ethyl)-2,4-dihydro-3H-1,2,4-triazol-3-one;
 5-(3-{[4-(4-chlorophenyl)piperidin-1-yl]sulfonyl}propyl)-2,4-dihydro-3H-1,2,4-triazol-3-one;
 5-(3-{[4-(4-chlorophenyl)piperazin-1-yl]sulfonyl}propyl)-2,4-dihydro-3H-1,2,4-triazol-3-one;
 and pharmaceutically acceptable salts and solvates thereof.

10. A process for the preparation of a compound of formula (I) or a pharmaceutically acceptable salt or solvate thereof as defined in claim 1 which comprises:

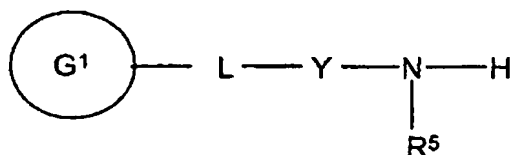
reaction of a compound of formula (II)



(II)

wherein R¹, R², R³, R⁴, X and m are as defined in Claim 1 and L¹ represents a leaving group, with a compound of formula (III)

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wherein G^1 , L, Y and R^5 are as defined in Claim 1;

5 and optionally thereafter forming a pharmaceutically acceptable salt or solvate.

11. A pharmaceutical composition comprising a compound of formula (I) or a pharmaceutically acceptable salt or solvate thereof as claimed in any one of claims 1 to 9 in association with a pharmaceutically acceptable adjuvant, diluent or carrier.

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12. A process for the preparation of a pharmaceutical composition as claimed in claim 11 which comprises mixing a compound of formula (I) or a pharmaceutically acceptable salt or solvate thereof as defined in any one of claims 1 to 9 with a pharmaceutically acceptable adjuvant, diluent or carrier.

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13. A compound of formula (I) or a pharmaceutically acceptable salt or solvate thereof as claimed in any one of claims 1 to 9 for use in therapy.

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14. Use of a compound of formula (I) or a pharmaceutically acceptable salt or solvate thereof as claimed in any one of claims 1 to 9 in the manufacture of a medicament for use in the treatment of an obstructive airways disease.

15. Use according to claim 14, wherein the obstructive airways disease is asthma or chronic obstructive pulmonary disease.

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16. A method of treating a disease or condition mediated by MMP12 and/or MMP9 which comprises administering to a patient a therapeutically effective amount of a compound of formula (I) or a pharmaceutically acceptable salt or solvate thereof as claimed in any one of claims 1 to 9.

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17. A method of treating an obstructive airways disease which comprises administering to a patient a therapeutically effective amount of a compound of formula (I) or a pharmaceutically acceptable salt or solvate thereof as claimed in any one of claims 1 to 9.

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